



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:24 PM UTC

PDB ID : 6C6N / pdb_00006c6n
Title : Human squalene epoxidase (SQLE, squalene monooxygenase) structure with FAD and Cmpd-4"
Authors : Padyana, A.K.; Jin, L.
Deposited on : 2018-01-19
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

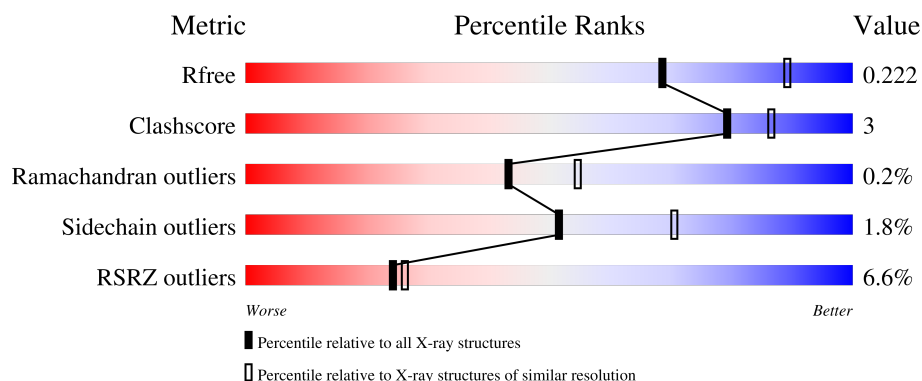
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	458	5% 92% 5% ..
1	B	458	7% 89% 9% .

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7638 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

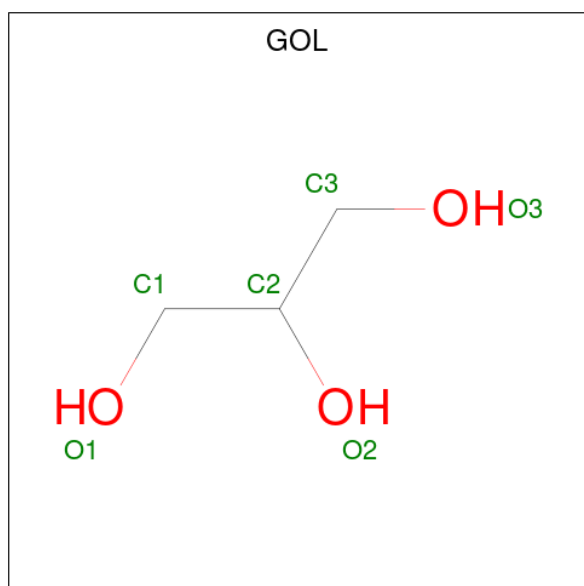
- Molecule 1 is a protein called Squalene monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	0	0
			3526	2280	594	637	15			
1	B	450	Total	C	N	O	S	0	1	0
			3532	2283	595	639	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	117	GLY	-	expression tag	UNP Q14534
B	117	GLY	-	expression tag	UNP Q14534

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



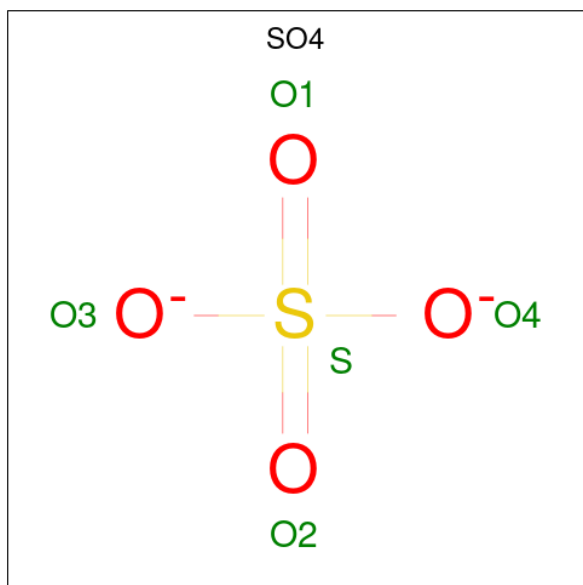
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		

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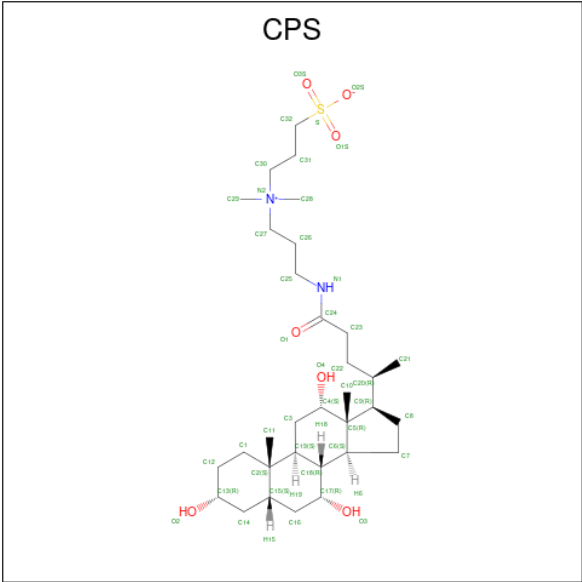
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



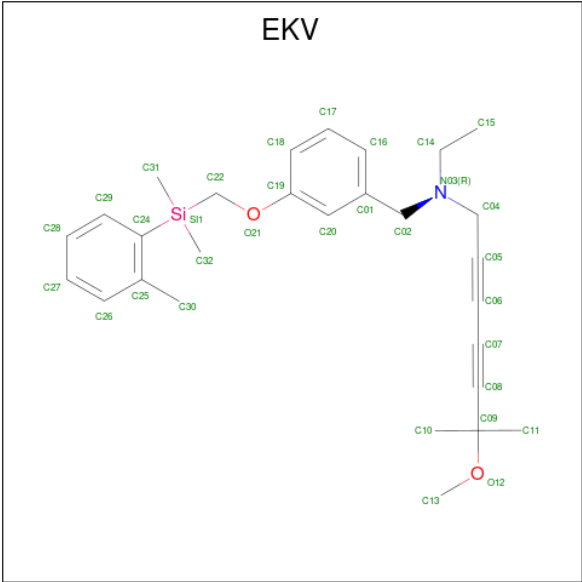
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 3-[(3-CHOLAMIDOPROPYL)DIMETHYLAMMONIO]-1-PROPANESULFONATE (CCD ID: CPS) (formula: C₃₂H₅₈N₂O₇S).



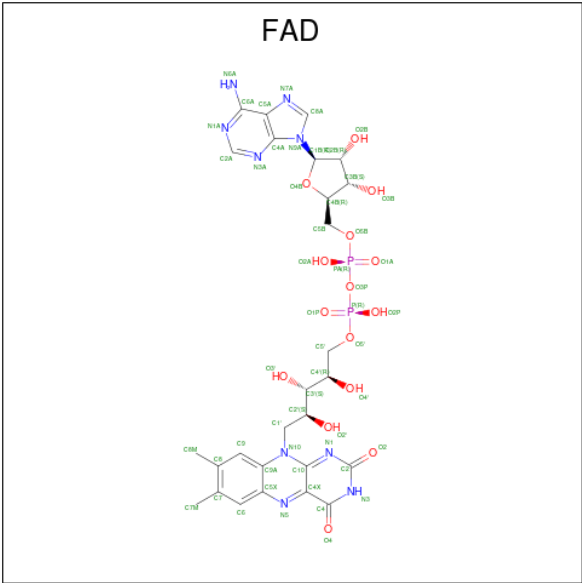
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O 31 26 1 4	0	0
4	A	1	Total C N O 30 25 1 4	0	0
4	A	1	Total C O 26 23 3	0	0
4	A	1	Total C N O 29 24 1 4	0	0
4	A	1	Total C O 25 22 3	0	0
4	B	1	Total C N O 30 25 1 4	0	0
4	B	1	Total C O 26 23 3	0	0

- Molecule 5 is N-[(3-{[dimethyl(2-methylphenyl)silyl]methoxy}phenyl)methyl]-N-ethyl-6-methoxy-6-methylhepta-2,4-diyne-1-amine (CCD ID: EKV) (formula: C₂₈H₃₇NO₂Si).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	Si	0	0
			32	28	1	2	1		
5	B	1	Total	C	N	O	Si	0	0
			32	28	1	2	1		

- Molecule 6 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
6	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

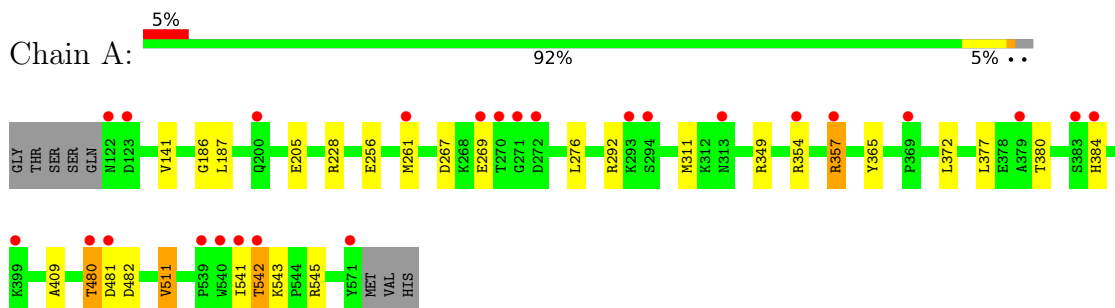
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	98	Total 98	O 98	0	0
7	B	81	Total 81	O 81	0	0

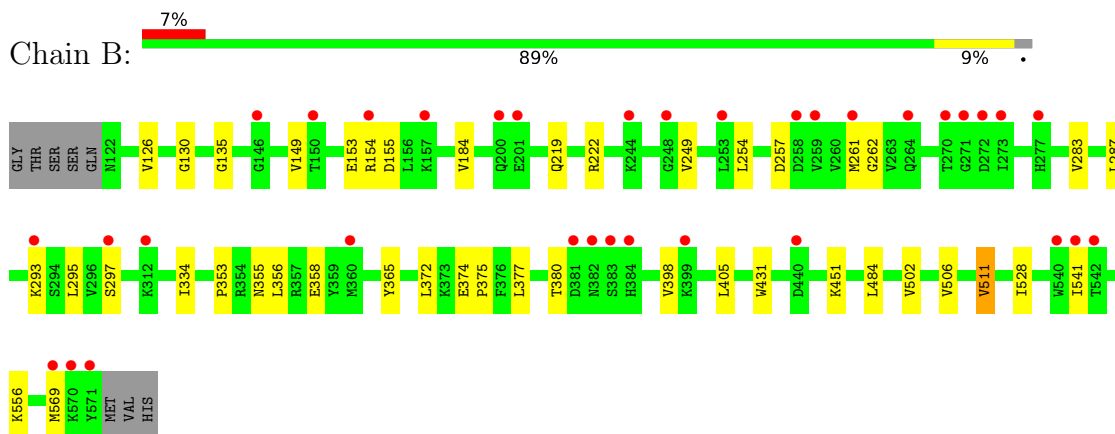
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Squalene monooxygenase



- Molecule 1: Squalene monooxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	126.96Å 126.96Å 166.12Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.16 – 2.30 37.16 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.8 (37.16-2.30) 98.8 (37.16-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.97 (at 2.29Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575)	Depositor
R, R_{free}	0.189 , 0.220 0.190 , 0.222	Depositor DCC
R_{free} test set	3351 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	43.3	Xtriage
Anisotropy	0.001	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7638	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.03% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FAD, GOL, EKV, CPS, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.13	0/3609	0.33	0/4887
1	B	0.12	0/3615	0.33	0/4895
All	All	0.13	0/7224	0.33	0/9782

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3526	0	3580	17	0
1	B	3532	0	3584	23	0
2	A	18	0	24	1	0
2	B	6	0	8	1	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	141	0	193	1	0
4	B	56	0	77	0	0
5	A	32	0	0	0	0
5	B	32	0	0	0	0
6	A	53	0	31	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	53	0	31	1	0
7	A	98	0	0	0	0
7	B	81	0	0	0	0
All	All	7638	0	7528	40	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (40) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:ARG:HG3	1:B:249:VAL:HG22	1.88	0.55
1:B:126:VAL:HB	1:B:149:VAL:HG22	1.94	0.50
1:B:334:ILE:HD13	1:B:372:LEU:HD13	1.93	0.50
1:A:365:TYR:CE1	1:A:377:LEU:HG	2.48	0.48
1:B:295:LEU:HB3	1:B:398:VAL:HG13	1.96	0.48
1:B:219:GLN:HE21	2:B:601:GOL:H32	1.78	0.47
1:A:481:ASP:OD1	1:A:482:ASP:N	2.47	0.47
1:B:431:TRP:CH2	1:B:451:LYS:HG3	2.50	0.47
1:B:484:LEU:HG	1:B:556:LYS:HG2	1.96	0.47
1:A:543:LYS:HG3	1:B:569:MET:HE2	1.96	0.47
1:B:374:GLU:HB3	1:B:375:PRO:HD3	1.96	0.47
1:B:431:TRP:HH2	1:B:451:LYS:HG3	1.80	0.47
1:B:283:VAL:HB	1:B:405:LEU:HD12	1.97	0.47
1:A:256:GLU:HG3	1:A:261:MET:HG3	1.96	0.46
1:A:267:ASP:OD1	1:A:269:GLU:HG2	2.15	0.46
1:A:186:GLY:HA2	2:A:602:GOL:H32	1.98	0.46
1:A:187:LEU:HD22	1:A:228:ARG:HB3	1.97	0.46
1:A:377:LEU:HD23	1:A:377:LEU:HA	1.82	0.46
1:A:542:THR:O	1:A:545:ARG:HB2	2.15	0.46
1:B:184:VAL:HG11	1:B:222:ARG:HG3	1.97	0.46
1:A:269:GLU:OE2	1:A:269:GLU:N	2.46	0.46
1:A:349:ARG:HH12	1:A:480:THR:HG21	1.81	0.46
1:B:153:GLU:HG3	1:B:155:ASP:H	1.82	0.45
1:B:377:LEU:O	1:B:380:THR:OG1	2.33	0.44
1:A:256:GLU:HB2	1:A:261:MET:HE3	1.99	0.44
1:A:311:MET:HE1	1:A:372:LEU:HD22	2.00	0.43
1:A:357:ARG:HE	1:A:357:ARG:HB3	1.61	0.43
1:B:502:VAL:O	1:B:506:VAL:HG23	2.18	0.43
1:B:254:LEU:N	1:B:262:GLY:O	2.51	0.42
4:A:605:CPS:H11A	1:B:528:ILE:HG13	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:355:ASN:CG	1:B:358:GLU:HB2	2.45	0.42
6:B:606:FAD:H9	6:B:606:FAD:H1'1	1.86	0.42
1:B:365:TYR:CE1	1:B:377:LEU:HG	2.55	0.41
1:B:293:LYS:HA	1:B:293:LYS:HD3	1.92	0.41
1:A:292:ARG:HH21	1:A:409:ALA:HA	1.85	0.41
1:A:276:LEU:HD23	1:A:276:LEU:HA	1.87	0.41
1:A:377:LEU:O	1:A:380:THR:OG1	2.36	0.41
1:B:287:LEU:H	1:B:287:LEU:HD23	1.86	0.41
1:B:130:GLY:O	1:B:135:GLY:HA3	2.21	0.40
1:B:353:PRO:HG2	1:B:356:LEU:HD23	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	448/458 (98%)	435 (97%)	12 (3%)	1 (0%)	43	55
1	B	449/458 (98%)	438 (98%)	10 (2%)	1 (0%)	43	55
All	All	897/916 (98%)	873 (97%)	22 (2%)	2 (0%)	43	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	511	VAL
1	B	511	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	384/391 (98%)	375 (98%)	9 (2%)	44	63
1	B	385/391 (98%)	380 (99%)	5 (1%)	61	77
All	All	769/782 (98%)	755 (98%)	14 (2%)	51	70

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	141	VAL
1	A	205	GLU
1	A	354	ARG
1	A	357	ARG
1	A	384	HIS
1	A	480	THR
1	A	511	VAL
1	A	541	ILE
1	A	542	THR
1	B	257	ASP
1	B	261	MET
1	B	297	SER
1	B	511	VAL
1	B	541	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	298	ASN
1	A	320	ASN
1	A	371	HIS
1	A	486	GLN
1	B	219	GLN
1	B	226	HIS
1	B	264	GLN
1	B	298	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

17 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	CPS	B	604	-	29,29,45	0.19	0	47,47,70	0.41	0
4	CPS	A	609	-	28,28,45	0.20	0	46,46,70	0.41	0
3	SO4	A	604	-	4,4,4	0.22	0	6,6,6	0.07	0
3	SO4	B	602	-	4,4,4	0.25	0	6,6,6	0.08	0
4	CPS	A	608	-	32,32,45	0.15	0	51,51,70	0.35	0
2	GOL	A	601	-	5,5,5	0.37	0	5,5,5	0.31	0
4	CPS	B	603	-	33,33,45	0.16	0	52,52,70	0.37	0
2	GOL	A	603	-	5,5,5	0.37	0	5,5,5	0.33	0
4	CPS	A	607	-	29,29,45	0.19	0	47,47,70	0.37	0
5	EKV	A	610	-	30,33,33	2.50	8 (26%)	33,45,45	1.06	3 (9%)
5	EKV	B	605	-	30,33,33	2.50	9 (30%)	33,45,45	0.90	1 (3%)
2	GOL	B	601	-	5,5,5	0.37	0	5,5,5	0.29	0
4	CPS	A	605	-	34,34,45	0.19	0	53,53,70	0.37	0
6	FAD	B	606	-	58,58,58	0.29	0	85,89,89	0.32	0
6	FAD	A	611	-	58,58,58	0.28	0	85,89,89	0.33	0
2	GOL	A	602	-	5,5,5	0.37	0	5,5,5	0.26	0
4	CPS	A	606	-	33,33,45	0.18	0	52,52,70	0.38	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	CPS	B	604	-	-	0/6/71/90	0/4/4/4
4	CPS	A	609	-	-	0/4/69/90	0/4/4/4
4	CPS	A	608	-	-	1/9/74/90	0/4/4/4
2	GOL	A	601	-	-	2/4/4/4	-
4	CPS	B	603	-	-	3/11/76/90	0/4/4/4
2	GOL	A	603	-	-	2/4/4/4	-
4	CPS	A	607	-	-	0/6/71/90	0/4/4/4
5	EKV	A	610	-	-	5/27/30/30	0/2/2/2
5	EKV	B	605	-	-	2/27/30/30	0/2/2/2
2	GOL	B	601	-	-	2/4/4/4	-
4	CPS	A	605	-	-	0/12/77/90	0/4/4/4
6	FAD	B	606	-	-	2/34/50/50	0/6/6/6
6	FAD	A	611	-	-	2/34/50/50	0/6/6/6
2	GOL	A	602	-	-	2/4/4/4	-
4	CPS	A	606	-	-	2/11/76/90	0/4/4/4

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	610	EKV	C09-C08	8.16	1.55	1.48
5	B	605	EKV	C09-C08	8.04	1.54	1.48
5	A	610	EKV	C07-C06	7.86	1.54	1.37
5	B	605	EKV	C07-C06	7.84	1.54	1.37
5	B	605	EKV	C04-C05	3.63	1.55	1.47
5	A	610	EKV	C04-C05	3.63	1.55	1.47
5	B	605	EKV	O21-C19	2.89	1.44	1.37
5	A	610	EKV	O21-C19	2.82	1.44	1.37
5	A	610	EKV	C02-C01	2.47	1.55	1.51
5	B	605	EKV	C02-C01	2.40	1.55	1.51
5	B	605	EKV	SI1-C24	2.30	1.93	1.88
5	A	610	EKV	SI1-C24	2.27	1.93	1.88
5	B	605	EKV	C04-N03	2.19	1.49	1.46
5	B	605	EKV	C10-C09	2.15	1.55	1.52
5	A	610	EKV	SI1-C32	2.15	1.91	1.86
5	B	605	EKV	SI1-C32	2.09	1.91	1.86
5	A	610	EKV	C10-C09	2.02	1.55	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	610	EKV	C10-C09-C08	-3.60	105.09	109.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	605	EKV	C29-C24-C25	2.49	120.03	117.86
5	A	610	EKV	C29-C24-C25	2.48	120.02	117.86
5	A	610	EKV	C05-C04-N03	-2.02	109.85	114.08

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	GOL	C1-C2-C3-O3
2	B	601	GOL	O1-C1-C2-C3
5	A	610	EKV	N03-C04-C05-C06
5	A	610	EKV	C10-C09-O12-C13
5	A	610	EKV	C08-C09-O12-C13
5	A	610	EKV	C11-C09-O12-C13
2	A	602	GOL	O1-C1-C2-C3
2	A	603	GOL	O1-C1-C2-C3
2	A	601	GOL	O2-C2-C3-O3
6	B	606	FAD	O4B-C4B-C5B-O5B
6	B	606	FAD	C3B-C4B-C5B-O5B
2	A	603	GOL	O1-C1-C2-O2
2	B	601	GOL	O1-C1-C2-O2
6	A	611	FAD	O4B-C4B-C5B-O5B
4	A	608	CPS	C20-C22-C23-C24
4	A	606	CPS	C22-C23-C24-O1
4	B	603	CPS	C20-C22-C23-C24
4	A	606	CPS	C22-C23-C24-N1
5	B	605	EKV	C29-C24-SI1-C22
6	A	611	FAD	C3B-C4B-C5B-O5B
5	B	605	EKV	C15-C14-N03-C04
5	A	610	EKV	C29-C24-SI1-C22
4	B	603	CPS	C22-C23-C24-O1
4	B	603	CPS	C22-C23-C24-N1
2	A	602	GOL	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 4 short contacts:

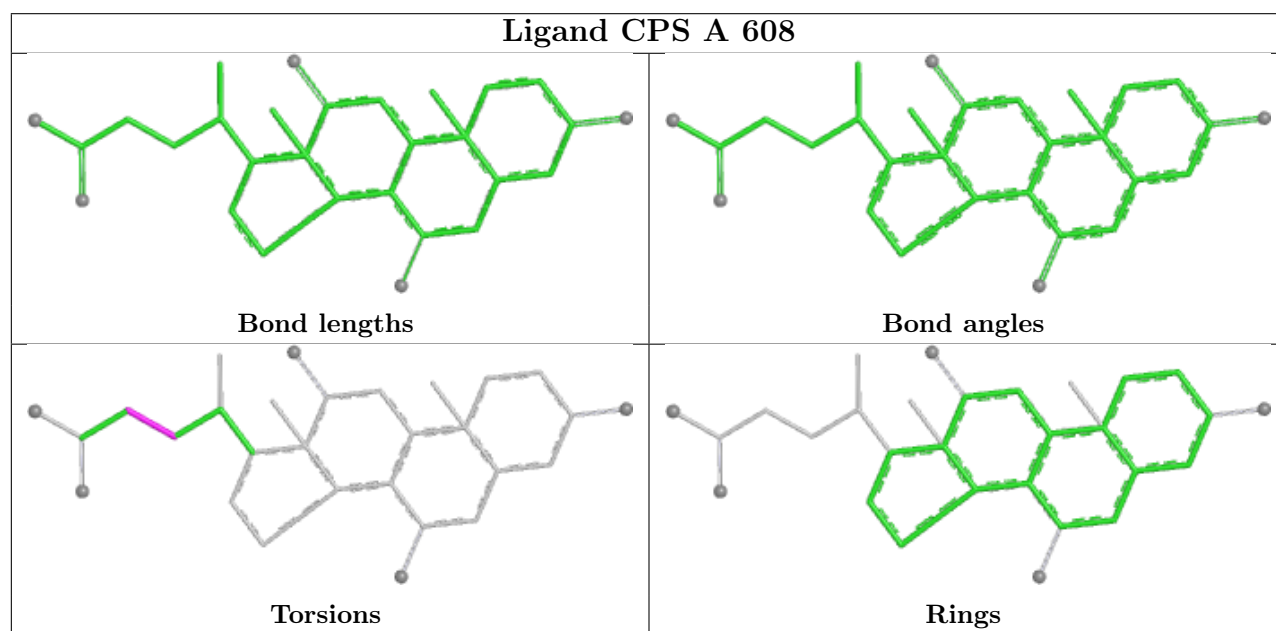
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	GOL	1	0
4	A	605	CPS	1	0
6	B	606	FAD	1	0

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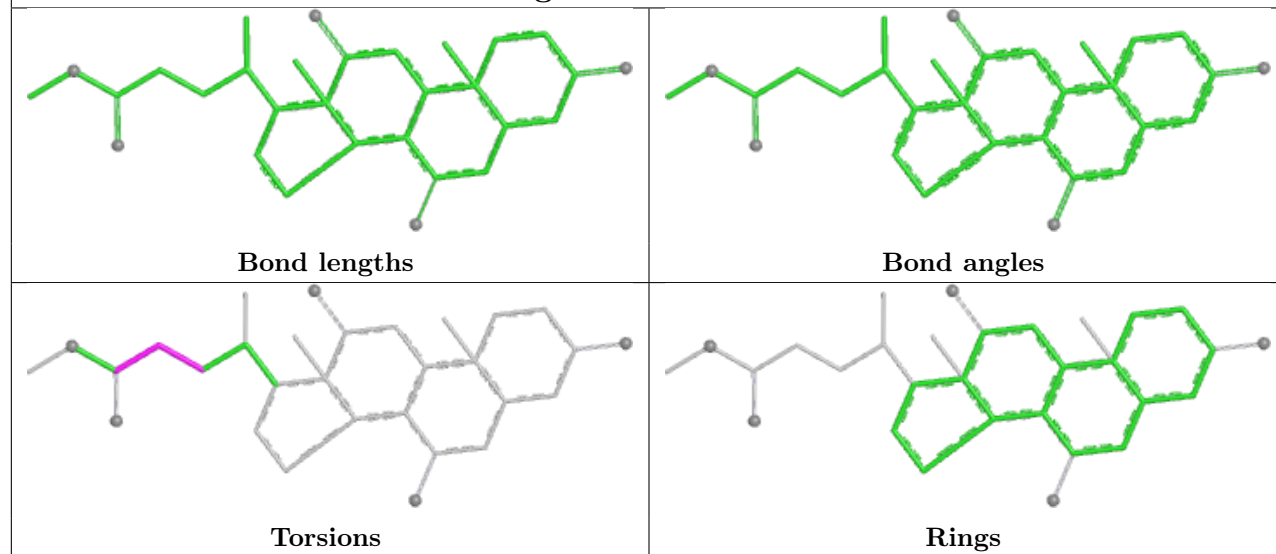
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	602	GOL	1	0

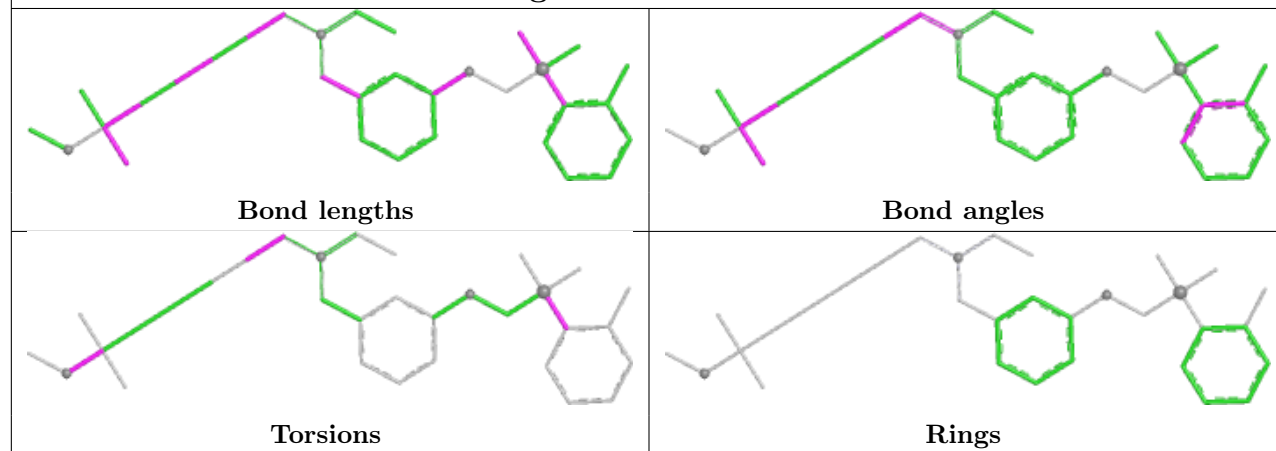
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



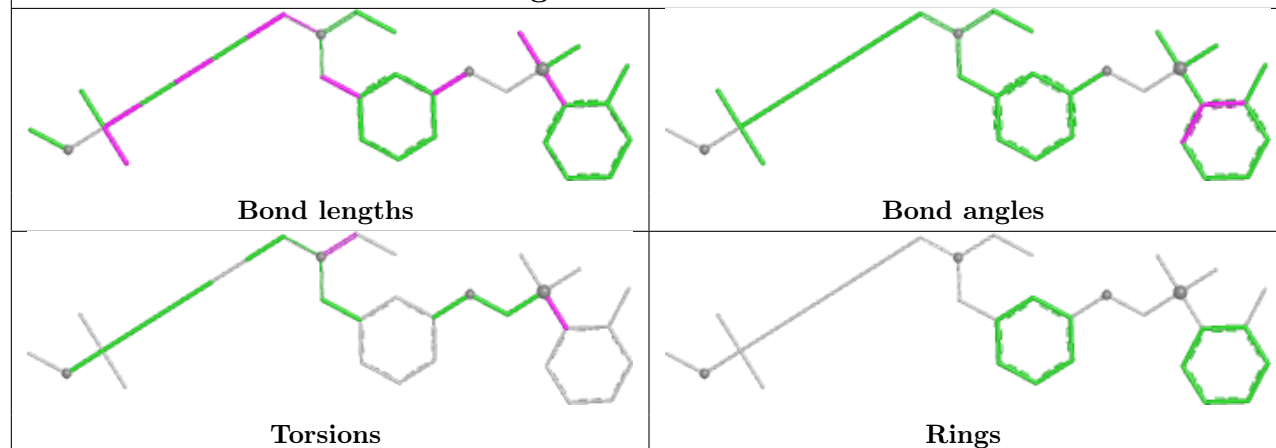
Ligand CPS B 603

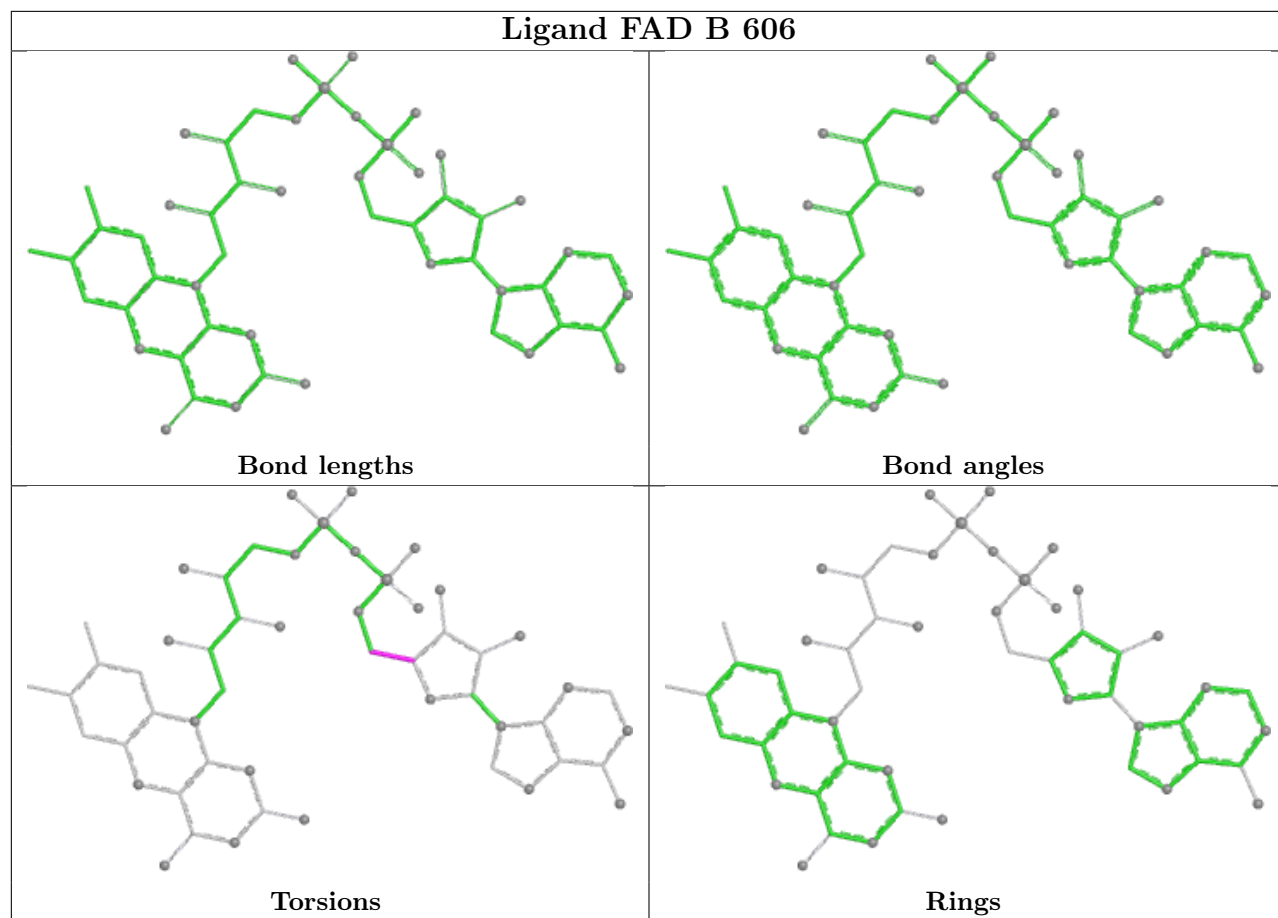


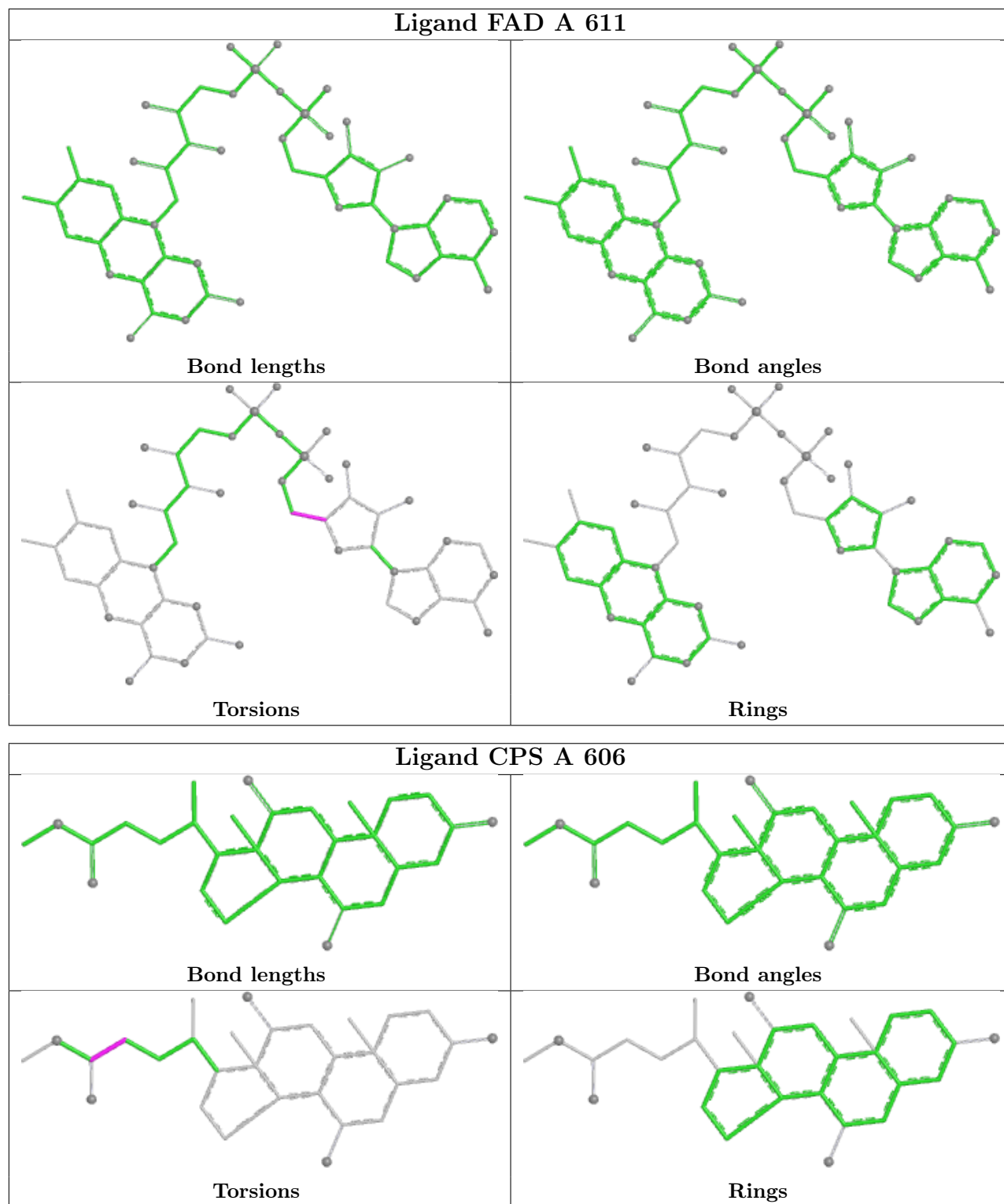
Ligand EKV A 610



Ligand EKV B 605







5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	450/458 (98%)	0.34	25 (5%) 30 32	28, 48, 88, 115	0
1	B	450/458 (98%)	0.49	34 (7%) 20 21	28, 50, 83, 109	1 (0%)
All	All	900/916 (98%)	0.42	59 (6%) 24 26	28, 49, 85, 115	1 (0%)

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	541	ILE	4.4
1	A	540	TRP	4.4
1	A	271	GLY	4.3
1	B	571	TYR	4.3
1	A	294	SER	4.2
1	B	272	ASP	4.0
1	B	259	VAL	3.8
1	B	541	ILE	3.8
1	B	570	LYS	3.7
1	A	313	ASN	3.6
1	B	277	HIS	3.6
1	A	571	TYR	3.6
1	B	293	LYS	3.5
1	B	382	ASN	3.4
1	A	269	GLU	3.1
1	A	270	THR	3.1
1	A	542	THR	3.0
1	B	271	GLY	3.0
1	B	381	ASP	3.0
1	B	261	MET	2.9
1	A	399	LYS	2.9
1	A	200	GLN	2.9
1	B	273	ILE	2.9
1	A	480	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	354	ARG	2.8
1	B	540	TRP	2.7
1	B	154	ARG	2.7
1	B	297	SER	2.6
1	B	360	MET	2.5
1	B	312	LYS	2.5
1	A	384	HIS	2.5
1	B	248	GLY	2.5
1	A	383	SER	2.5
1	B	440	ASP	2.4
1	B	258	ASP	2.3
1	A	369	PRO	2.3
1	B	146	GLY	2.3
1	B	384	HIS	2.3
1	A	272	ASP	2.3
1	B	542	THR	2.3
1	A	293	LYS	2.3
1	A	379	ALA	2.2
1	B	201	GLU	2.2
1	A	122	ASN	2.2
1	A	123	ASP	2.2
1	A	481	ASP	2.2
1	A	539	PRO	2.2
1	B	270	THR	2.2
1	B	253	LEU	2.1
1	B	200	GLN	2.1
1	A	357	ARG	2.1
1	B	150	THR	2.1
1	B	383	SER	2.1
1	A	261	MET	2.1
1	B	264	GLN	2.0
1	B	157	LYS	2.0
1	B	244	LYS	2.0
1	B	399	LYS	2.0
1	B	569	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

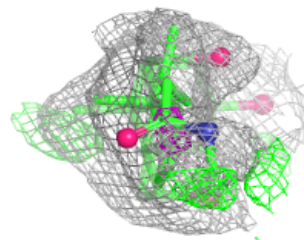
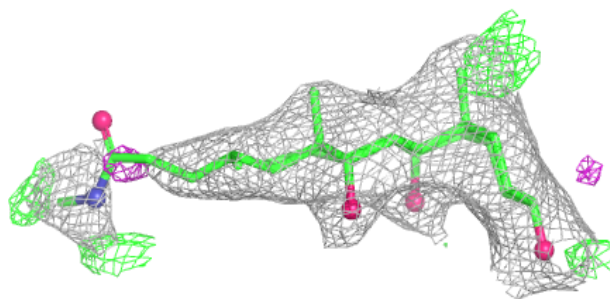
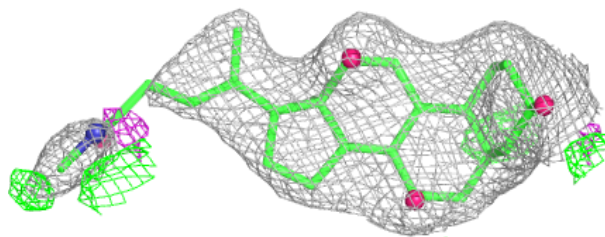
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GOL	A	603	6/6	0.73	0.24	82,82,92,92	0
2	GOL	B	601	6/6	0.77	0.18	60,70,76,81	0
2	GOL	A	602	6/6	0.78	0.20	59,75,83,84	0
2	GOL	A	601	6/6	0.82	0.15	54,61,66,67	0
4	CPS	A	607	26/42	0.86	0.15	64,78,87,90	0
4	CPS	B	604	26/42	0.86	0.17	64,75,81,83	0
3	SO4	A	604	5/5	0.88	0.17	67,69,78,81	5
4	CPS	B	603	30/42	0.89	0.14	61,74,95,105	0
3	SO4	B	602	5/5	0.90	0.13	78,81,103,107	0
4	CPS	A	608	29/42	0.91	0.14	60,68,96,103	0
4	CPS	A	609	25/42	0.92	0.14	39,54,71,75	0
4	CPS	A	605	31/42	0.93	0.11	28,38,83,109	0
4	CPS	A	606	30/42	0.94	0.11	31,41,79,88	0
6	FAD	A	611	53/53	0.95	0.09	36,43,54,58	0
6	FAD	B	606	53/53	0.95	0.09	35,44,67,69	0
5	EKV	A	610	32/32	0.96	0.08	31,41,56,61	0
5	EKV	B	605	32/32	0.96	0.08	27,41,51,60	0

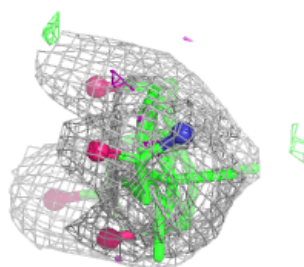
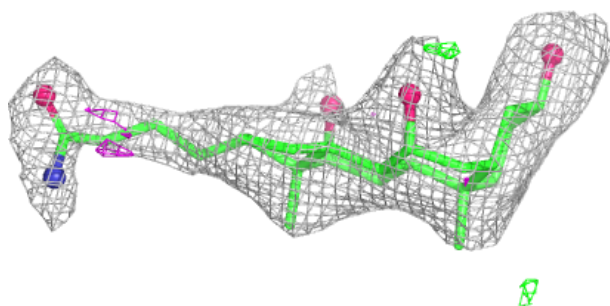
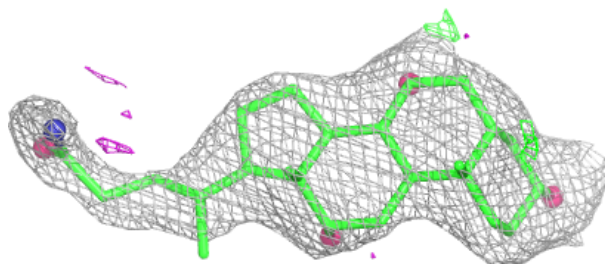
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around CPS B 603:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

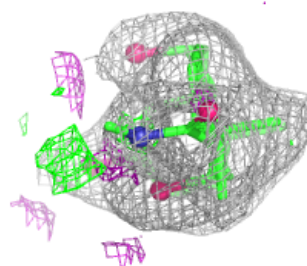
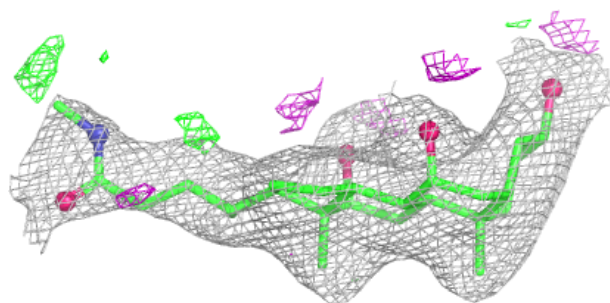
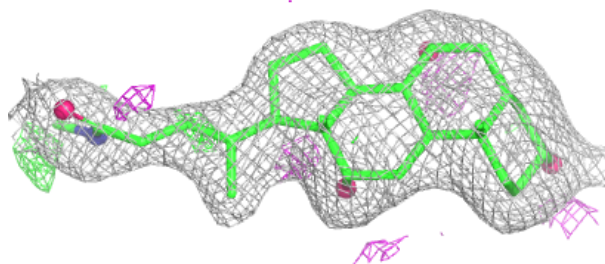
**Electron density around CPS A 608:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

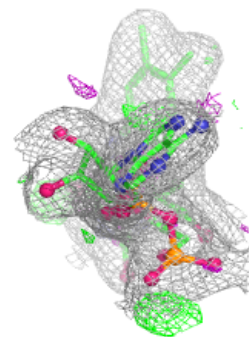
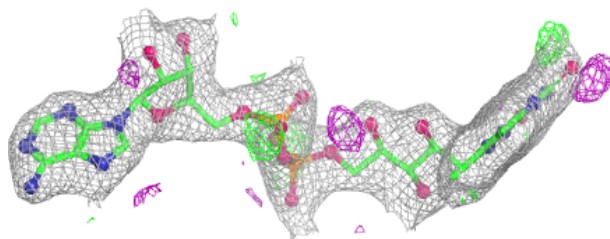
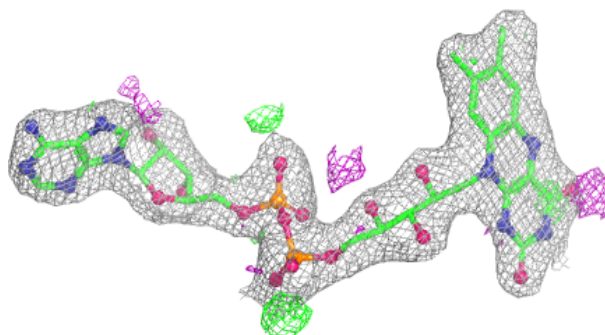


Electron density around CPS A 606:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

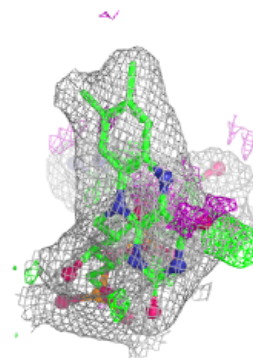
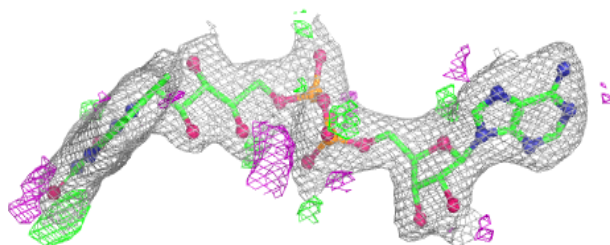
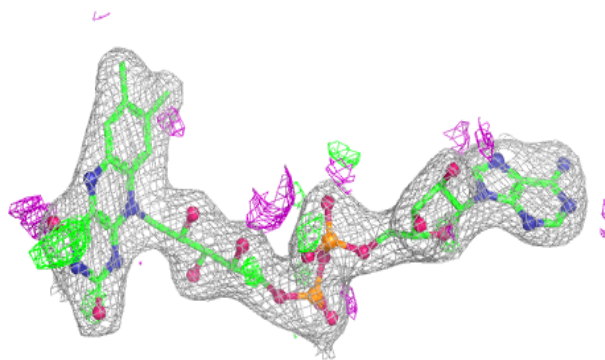
**Electron density around FAD A 611:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

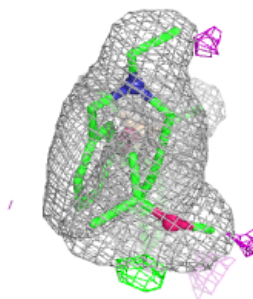
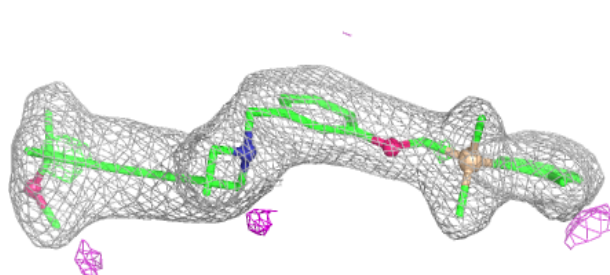
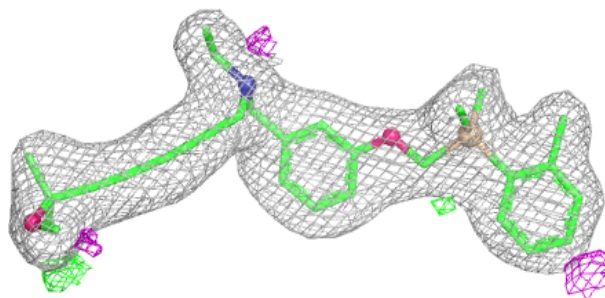


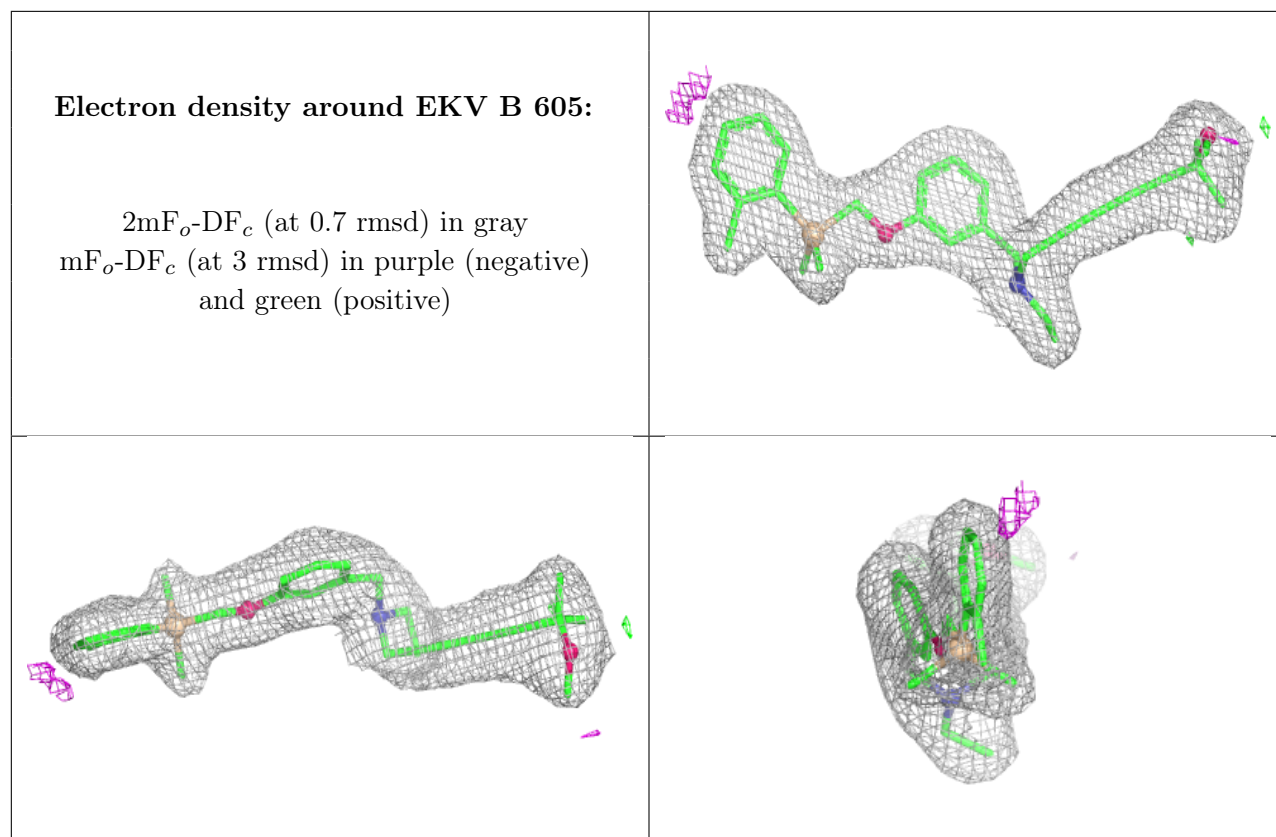
Electron density around FAD B 606:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around EKV A 610:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.